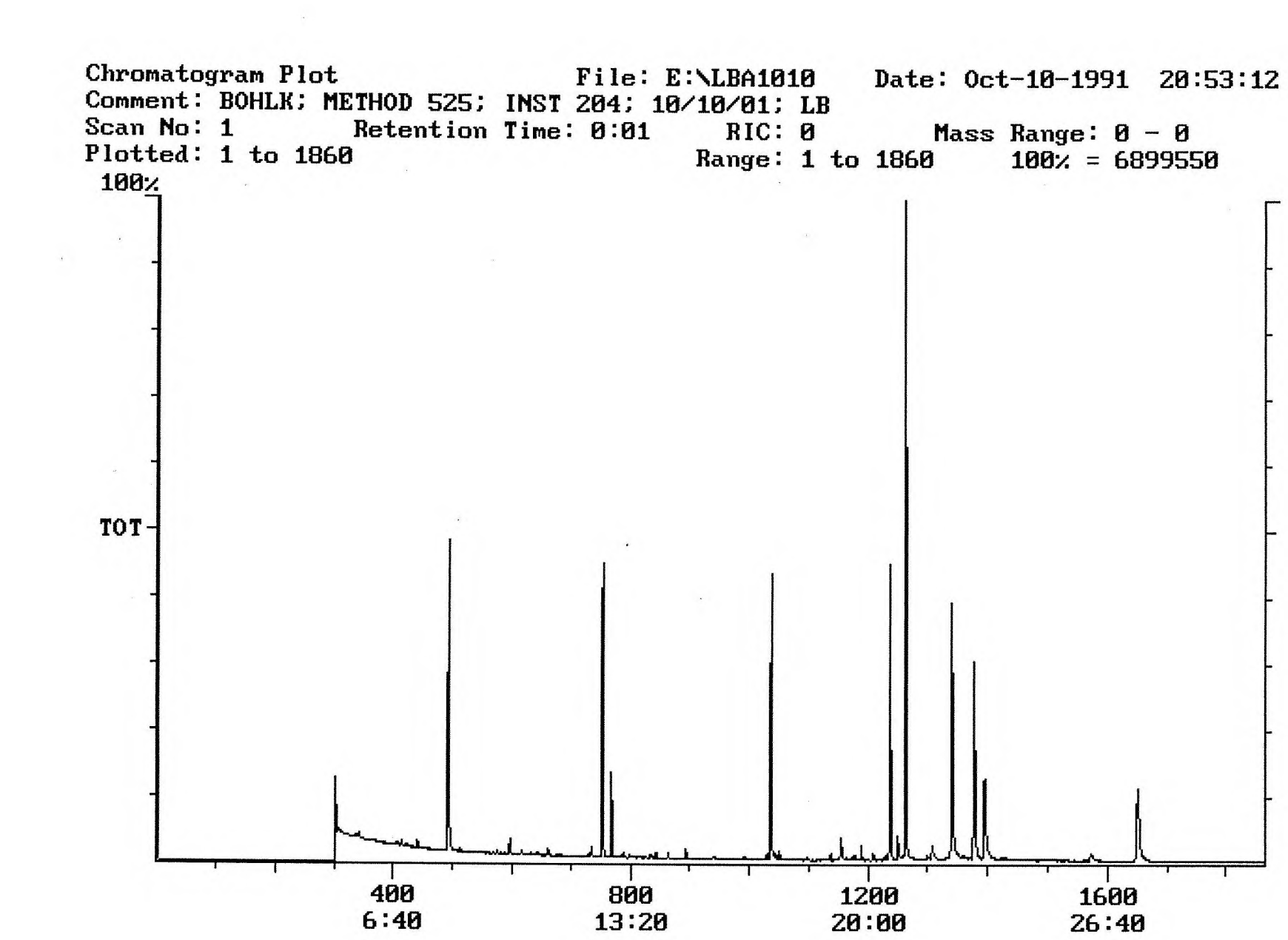
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 10/15/2001 Time: 11:06:30

Sample: AD23343
Analysis: $B1525 Organic Chemicals - 525.2; lab blank
Units: ug
Started: 10/4/01 11:00 Ended: 10/10/01 11:00 Analyst: TB
Result source: manual entry
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Hexachlorocyclopentadiene		< MDL		0.10
2	Hexachlorobenzene		< MDL		0.10
3	Simazine		< MDL		0.07
4	Atrazine		< MDL		0.10
5	Alachlor		< MDL		0.20
6	bis(2-Ethylhexyl)adipate		< MDL		0.60
7	bis(2-Ethylhexyl)phthalate	USPEC	< MDL		0.10
8	Benzo(a)pyrene		< MDL		0.20
9	EPTC		< MDL		1.0
10	2,6-Dinitrotoluene		< MDL		2.0
11	2,4-Dinitrotoluene		< MDL		2.0
12	Molinate		< MDL		0.90
13	Terbacil		< MDL		2.0
14	Acetochlor		< MDL		2.0
15	Metribuzin		< MDL		0.15
16	Metolachlor		< MDL		0.75
17	Butachlor		< MDL		0.40
18	4,4-DDE		< MDL		0.80
19	Surr_1,3-Dimethyl-2-nitrobenzen		5.5		
20	Surr_Triphenyl phosphate		5.3		
21	Surr_Fluoranthene-d10	USPEC			



Integration Report Quanfile: LBA1010 Quan Entries: 16
Comment: BOHLK; METHOD 525; INST 204; 10/10/01; LB
Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:33	753	Auto	1,459,572	695,610
2	phenanthrene d-10 (IS)	17:16	1036	Auto	2,395,872	1,060,658
3	chrysene d-12(IS)	22:57	1377	Auto	1,537,341	610,093
4	p-terphenyl d-14(IS)	21:02	1262	Auto	3,095,138	2,233,461
5	acenaphthene d-10(SURR	12:33	753	Auto	1,459,572	695,610
6	phenanthrene d-10(SURR	17:16	1036	Auto	2,395,872	1,060,658
7	chrysene d-12 (SURR)	22:57	1377	Auto	1,537,341	610,093
8	1,3-dimethyl-2-nitrobe	8:12	492	Auto	928,259	508,189
9	triphenyl phosphate (S	22:20	1340	Auto	1,213,147	392,070
10	perylene d-12 (SURR)	27:31	1651	Auto	1,337,660	235,688
11	hexachlorobenzene	15:43	943	Auto	5,370	2,393
12	bis(2-ethylhexyl)adipa	22:20	1340	Auto	10,620	4,032
13	bis(2-ethylhexyl)phtha	23:15	1395	Auto	1,012,705	241,039
14	2,6-dinitrotoluene	12:12	732	Auto	18,915	8,359
15	2,4-dinitrotoluene	13:19	799	Auto	903	903
16	4,4'-dde	20:58	1258	Auto	5,405	3,199

Quantitation Report Quanfile: LBA1010 Quan Entries: 16
Comment: BOHLK; METHOD 525; INST 204; 10/10/01; LB
Sorted via: Entry Number ↑ (S) = Standard

Cal	Name of Compound	S	Scan#	R Time	Me	Calc Amt(A)	Units
1	acenaphthene d-10(IS)	I	753	12:33	BV	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1036	17:16	BB	5.000	UG/L
3	chrysene d-12(IS)	I	1377	22:57	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1262	21:02	BB	5.000	UG/L
5	acenaphthene d-10(SURR	A	753	12:33	BV	2.987	UG/L
6	phenanthrene d-10(SURR	A	1036	17:16	BB	3.488	UG/L
7	chrysene d-12 (SURR)	A	1377	22:57	BB	3.371	UG/L
8	1,3-dimethyl-2-nitrobe	A	492	8:12	BV	5.461	UG/L
9	triphenyl phosphate (S	A	1340	22:20	BB	5.339	UG/L
10	perylene d-12 (SURR)	A	1651	27:31	BB	7.361	UG/L
12	hexachlorobenzene	A	943	15:43	BB	0.029	UG/L
16	bis(2-ethylhexyl)adipa	A	1340	22:20	MM	0.046	UG/L
17	bis(2-ethylhexyl)phtha	A	1395	23:15	VB	2.174	UG/L
20	2,6-dinitrotoluene	A	732	12:12	BB	0.187	UG/L
21	2,4-dinitrotoluene	A	799	13:19	BB	0.008	UG/L
28	4,4'-dde	A	1258	20:58	BB	0.023	UG/L

reviewed IS/surr's
Checked manually for Imagine
at mass 201 N.P.